

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW110318\
 Data File : VW006589.D
 Acq On : 02 Nov 2018 05:54
 Operator : SY/AP
 Sample : J5737-01
 Misc : 6.12G/5ML/MSVOA W/SOIL
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 TP-A

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W110218S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.678	284	291	301	rVB2	3848	11331	1.26%	0.177%
2	4.916	480	494	510	rBV2	19472	65718	7.31%	1.024%
3	7.891	972	982	986	rBV	145281	340027	37.81%	5.300%
4	7.952	986	992	1002	rVB	262987	580857	64.59%	9.053%
5	8.311	1041	1051	1063	rBV	124909	280773	31.22%	4.376%
6	8.848	1126	1139	1151	rBV	391389	790771	87.94%	12.325%
7	10.329	1374	1382	1388	rBV	492288	856906	95.29%	13.356%
8	10.390	1388	1392	1399	rVB	31055	52734	5.86%	0.822%
9	11.408	1549	1559	1560	rBV6	3866	10233	1.14%	0.159%
10	11.469	1560	1569	1576	rVV8	3460	15870	1.76%	0.247%
11	11.634	1589	1596	1605	rVB	539099	899266	100.00%	14.016%
12	12.621	1751	1758	1768	rVB	354440	583868	64.93%	9.100%
13	12.920	1789	1807	1809	rBV5	7927	30095	3.35%	0.469%
14	13.359	1862	1879	1893	rBV2	26683	131106	14.58%	2.043%
15	13.487	1893	1900	1906	rVV4	6078	21167	2.35%	0.330%
16	13.566	1906	1913	1920	rVB	515115	830529	92.36%	12.945%
17	13.762	1941	1945	1952	rVB	6086	10520	1.17%	0.164%
18	13.908	1955	1969	1970	rBV5	8645	21508	2.39%	0.335%
19	13.993	1975	1983	1993	rVV6	12869	46953	5.22%	0.732%
20	14.085	1993	1998	2011	rVB4	5595	13095	1.46%	0.204%
21	14.335	2031	2039	2050	rVB3	9481	26738	2.97%	0.417%
22	14.572	2067	2078	2089	rBV3	7814	23611	2.63%	0.368%
23	14.987	2128	2146	2160	rBV2	147464	536813	59.69%	8.367%
24	15.121	2160	2168	2175	rVV4	6541	21271	2.37%	0.332%
25	15.371	2202	2209	2219	rVB3	9241	24431	2.72%	0.381%
26	15.554	2237	2239	2247	rVB7	6476	12179	1.35%	0.190%
27	15.688	2247	2261	2275	rBV2	38746	163466	18.18%	2.548%
28	15.993	2301	2311	2319	rBV6	4218	14186	1.58%	0.221%

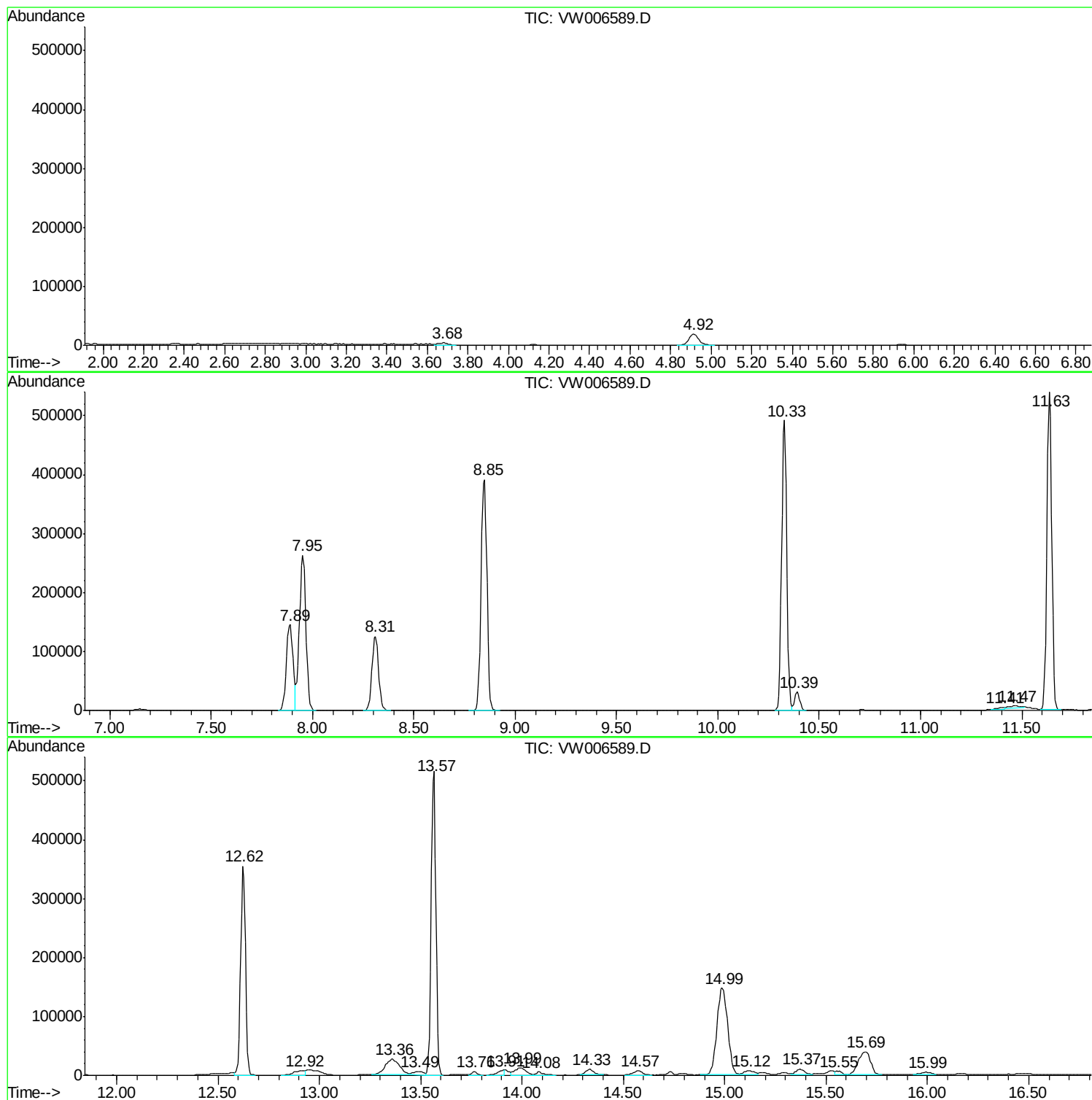
Sum of corrected areas: 6416022

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W110218S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
MSVOA_W
ClientSampled :
TP-A

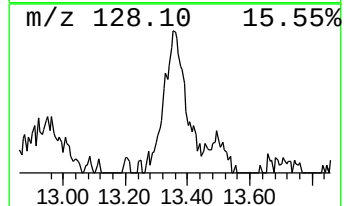
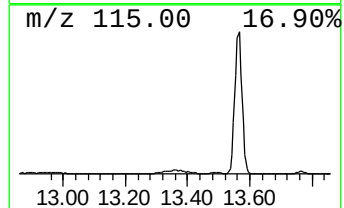
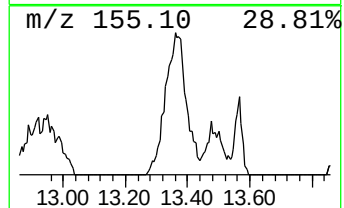
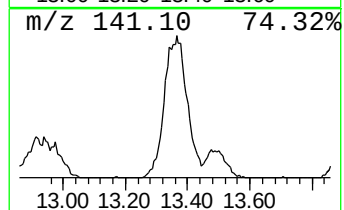
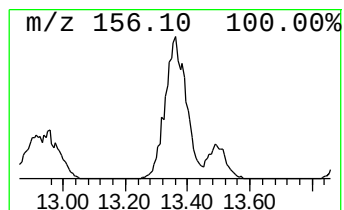
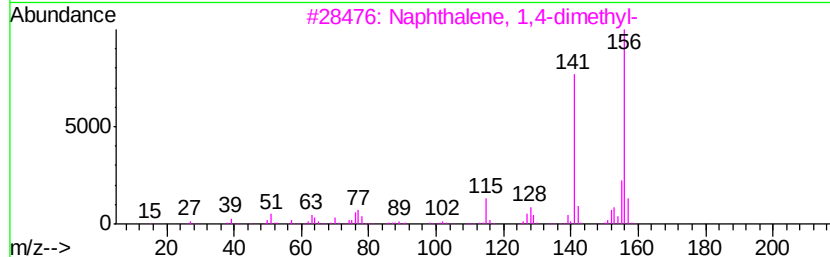
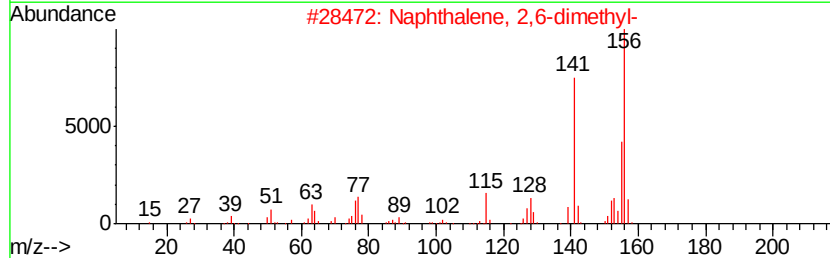
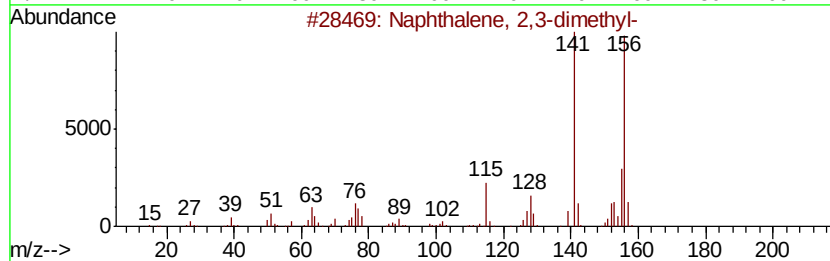
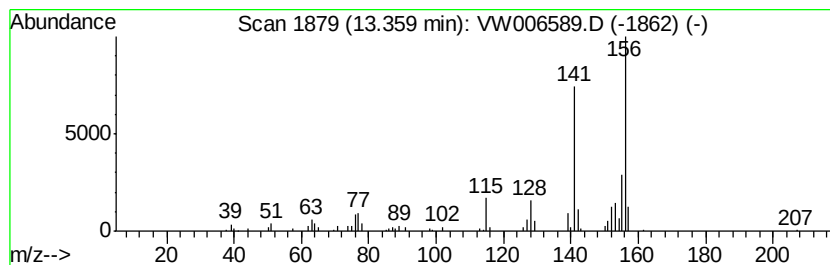
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W110218S.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 Naphthalene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	7.89 ug/l	131106	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



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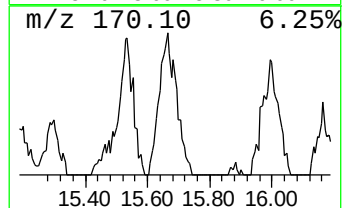
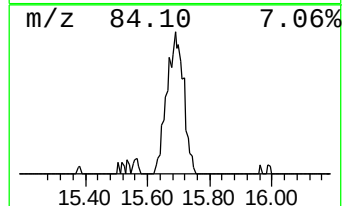
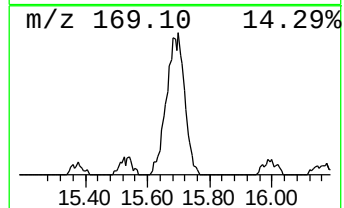
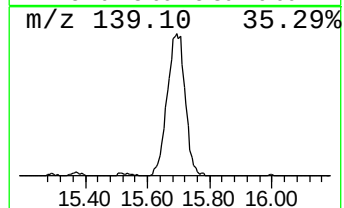
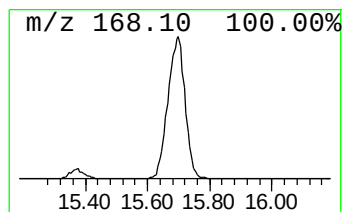
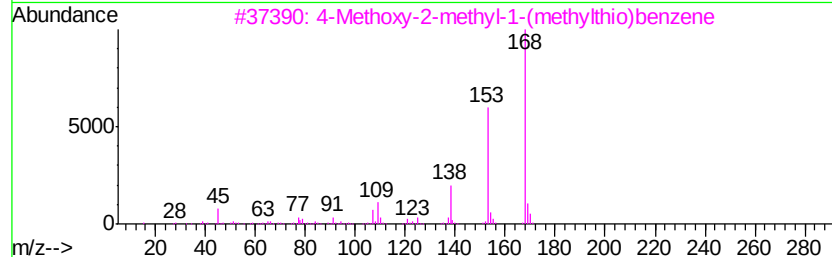
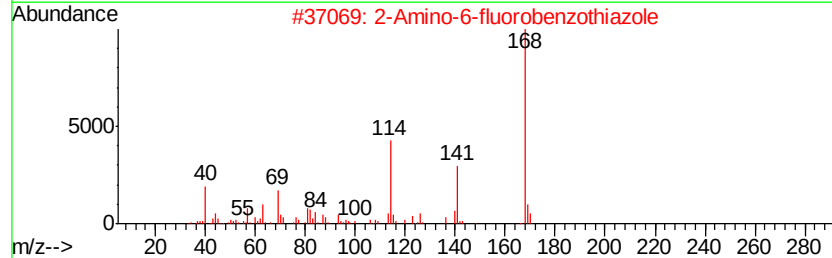
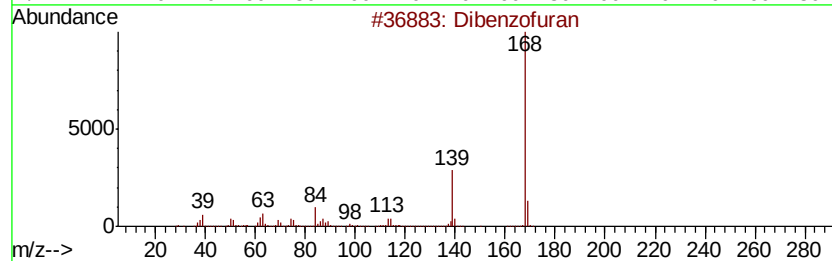
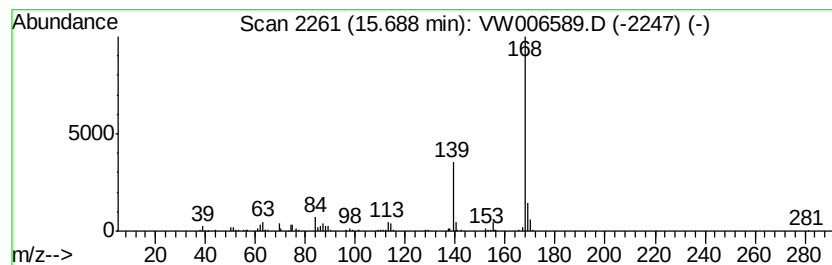
Instrument :
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Dibenzofuran Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.69	9.84 ug/l	163466	1,4-Dichlorobenzene-d4	13.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	90
2	2-Amino-6-fluorobenzothiazole	168	C7H5FN2S	000348-40-3	58
3	4-Methoxy-2-methyl-1-(methylthio)...	168	C9H12OS	022583-04-6	53
4	1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	53
5	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Naphthalene, 2,3-...	13.36	7.9	ug/l	131106	4	13.57	830529	50.0
Dibenzofuran	15.69	9.8	ug/l	163466	4	13.57	830529	50.0